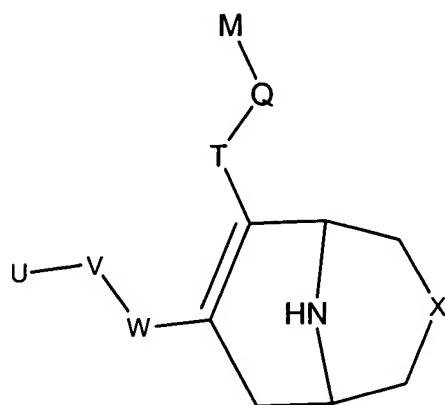


AMENDMENTS TO THE CLAIMS

The following list of claims will replace all prior claims in the application:

1. (Currently amended) A compound of formula I



Formula I

wherein

X represents -O-; -S-; -SO-; or -SO₂-;

W is a ~~six-membered, non-benzofused, phenyl or heteroaryl ring, substituted by V in meta or para position~~ non-benzofused phenyl or a six-membered, non-benzofused heteroaryl, wherein the phenyl and heteroaryl are substituted by V in meta or para position;

V represents a bond; -(CH₂)_r-; -A-(CH₂)_s-; -CH₂-A-(CH₂)_t-; -(CH₂)_s-A-; -(CH₂)₂-A-(CH₂)_u-; -A-(CH₂)_v-B-; -CH₂-CH₂-CH₂-A-CH₂-; -A-CH₂-CH₂-B-CH₂-; -CH₂-A-CH₂-CH₂-B-; -CH₂-CH₂-CH₂-A-CH₂-CH₂-; -CH₂-CH₂-CH₂-CH₂-A-CH₂-; -A-CH₂-CH₂-B-CH₂-CH₂-; -CH₂-A-CH₂-CH₂-B-CH₂-; -CH₂-A-CH₂-CH₂-CH₂-B-; -O-CH₂-CH(OCH₃)-CH₂-O-; -O-CH₂-CH(CH₃)-CH₂-O-; -O-CH₂-CH(CF₃)-CH₂-O-; -O-CH₂-C(CH₃)₂-CH₂-O-; -O-CH₂-C(CH₃)₂-O-; -O-C(CH₃)₂-CH₂-O-; -O-CH₂-CH(CH₃)-O-; -O-CH(CH₃)-CH₂-O-; -O-CH₂-C(CH₂CH₂)-O-; or -O-C(CH₂CH₂)-CH₂-O-;

A and B independently represent -O-; -S-; -SO-; or -SO₂-;

U represents aryl; or heteroaryl;

T represents -CONR¹-; -(CH₂)_pOCO-; -(CH₂)_pN(R¹)CO-; -(CH₂)_pN(R¹)SO₂-; or

-COO-;

Q represents lower alkylene; or lower alkenylene;

M represents hydrogen; cycloalkyl; aryl; heterocyclyl; or heteroaryl;

R¹ represents hydrogen; lower alkyl; lower alkenyl; lower alkynyl; cycloalkyl; aryl;
or cycloalkyl-lower alkyl;

p is the integer 1, 2, 3 or 4;

r is the integer 3, 4, 5, or 6;

s is the integer 2, 3, 4, or 5;

t is the integer 1, 2, 3, or 4;

u is the integer 1, 2, or 3; and

v is the integer 2, 3, or 4;

or optically pure enantiomers, racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, or the meso-form of the compound; or pharmaceutically acceptable salts, ~~solvent complexes or morphological forms~~ of the compound.

2. (Currently amended) The compound of formula I according to claim 1,
wherein

T represents -CONR¹-;

Q represents methylene; and

M represents aryl, or heteroaryl;

or optically pure enantiomers, racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, or the meso-form of the compound; or pharmaceutically acceptable salts, ~~solvent complexes or morphological forms~~ of the compound.

3. (Currently amended) The compound of formula I according to claim 1, wherein V represents -CH₂CH₂O-; -CH₂CH₂CH₂O-; or -OCH₂CH₂O-;

or optically pure enantiomers, racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, or the meso-form of the compound; or pharmaceutically acceptable salts, ~~solvent complexes or morphological forms~~ of the compound.

4. (Currently amended) The compound of formula I according to claim 1, wherein W represents a 1,4-disubstituted phenyl group;

or optically pure enantiomers, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, or the meso-form of the compound; or pharmaceutically acceptable salts, ~~solvent complexes or morphological forms~~ of the compound.

5. (Currently amended) The compound of formula I according to claim 1, wherein U is a mono-, di-, or trisubstituted phenyl or heteroaryl, whereby the substituents are selected from the group consisting of halogen, lower alkyl, lower alkoxy, and CF₃

or optically pure enantiomers, racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, or the meso-form of the compound; or pharmaceutically acceptable salts, ~~solvent complexes or morphological forms~~ of the compound.

6. (Currently amended) The compound according to claim 1 selected from the group consisting of:

[[*(rac.)*]](1*R**, 5*S**)-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3-oxa-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxy-2-methyl-benzyl)amide,

[[*(rac.)*]](1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3,3-dioxo-3λ⁶-thia-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide,

[[*(rac.)*]](1*R*^{*}, 3*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3-oxo-3λ⁴-thia-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxy-2-methylbenzyl)amide,

[[*(rac.)*]](1*R*^{*}, 3*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3-oxo-3λ⁴-thia-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methoxy-3-methylpyridin-4-ylmethyl)amide,

[[*(rac.)*]](1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3-oxa-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3-hydroxy-propoxy)-3-methylpyridin-4-ylmethyl]amide, and

[[*(rac.)*]](1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3,3-dioxo-3λ⁶-thia-9-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3-hydroxypropoxy)-3-methylpyridin-4-ylmethyl]amide.

7. (Previously presented) A pharmaceutical composition comprising at least one compound of claim 1 and a carrier and/or an adjuvant.

8-10. (Cancelled).